



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:09 PM GMT

PDB ID : 4HDH
Title : Crystal Structure of viral RdRp in complex with ATP
Authors : Surana, P.; Nair, D.T.
Deposited on : 2012-10-02
Resolution : 2.28 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

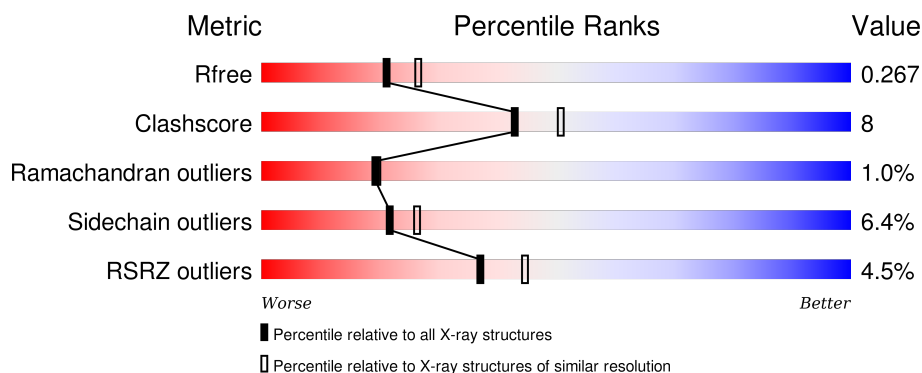
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


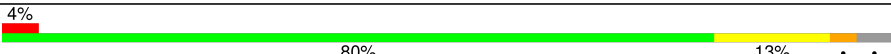
The reported resolution of this entry is 2.28 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	5193 (2.30-2.26)
Clashscore	102246	5929 (2.30-2.26)
Ramachandran outliers	100387	5851 (2.30-2.26)
Sidechain outliers	100360	5850 (2.30-2.26)
RSRZ outliers	91569	5204 (2.30-2.26)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	639	
1	B	639	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ATP	A	1001	-	-	-	X
2	ATP	B	1001	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10309 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Polypeptide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	613	Total	C	N	O	S	0	0	0
			4920	3104	886	899	31			
1	B	613	Total	C	N	O	S	0	0	0
			4904	3095	879	899	31			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	267	GLY	-	EXPRESSION TAG	UNP G3LHD9
A	268	PRO	-	EXPRESSION TAG	UNP G3LHD9
A	269	LEU	-	EXPRESSION TAG	UNP G3LHD9
A	270	GLY	-	EXPRESSION TAG	UNP G3LHD9
A	271	SER	-	EXPRESSION TAG	UNP G3LHD9
A	373	ARG	LYS	CONFLICT	UNP G3LHD9
A	429	ASN	ASP	CONFLICT	UNP G3LHD9
A	836	ALA	THR	CONFLICT	UNP G3LHD9
B	267	GLY	-	EXPRESSION TAG	UNP G3LHD9
B	268	PRO	-	EXPRESSION TAG	UNP G3LHD9
B	269	LEU	-	EXPRESSION TAG	UNP G3LHD9
B	270	GLY	-	EXPRESSION TAG	UNP G3LHD9
B	271	SER	-	EXPRESSION TAG	UNP G3LHD9
B	373	ARG	LYS	CONFLICT	UNP G3LHD9
B	429	ASN	ASP	CONFLICT	UNP G3LHD9
B	836	ALA	THR	CONFLICT	UNP G3LHD9

- Molecule 2 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	B	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total	Zn	0	0
			2	2		
3	A	2	Total	Zn	0	0
			2	2		

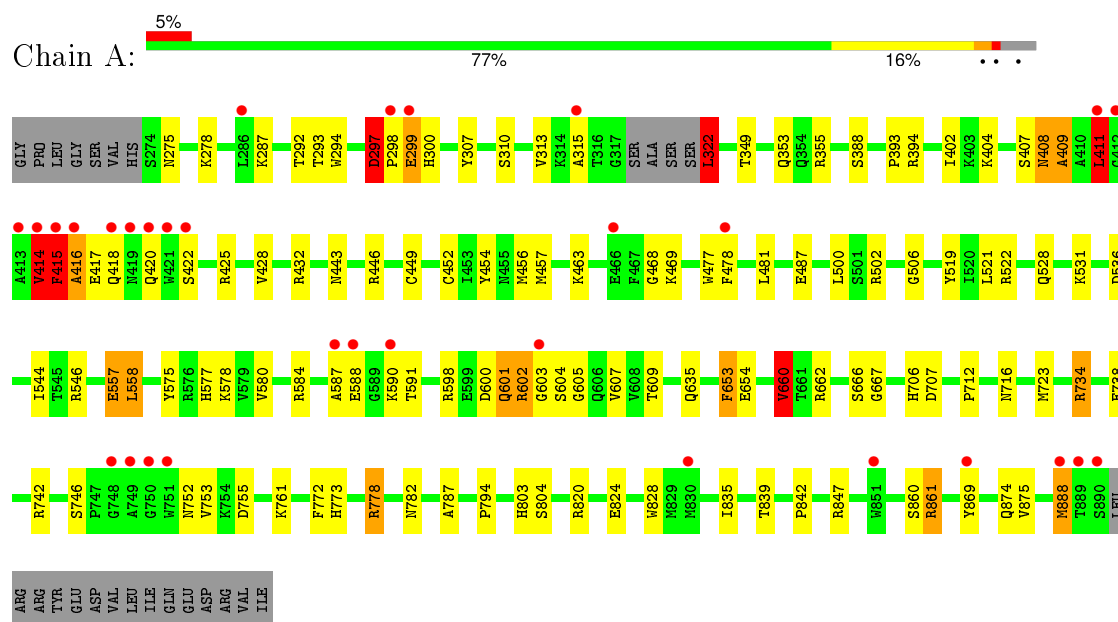
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	200	Total	O	0	0
			200	200		
4	B	219	Total	O	0	0
			219	219		

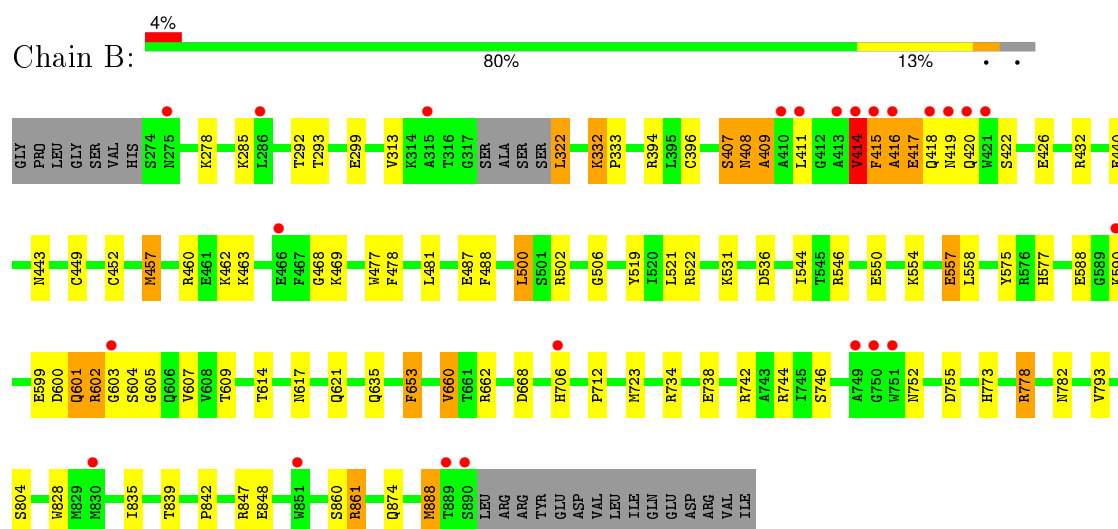
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Polyprotein



• Molecule 1: Polyprotein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	144.84Å 86.51Å 112.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	144.84 – 2.28 44.92 – 2.28	Depositor EDS
% Data completeness (in resolution range)	99.0 (144.84-2.28) 99.1 (44.92-2.28)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.06 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.225 , 0.272 0.220 , 0.267	Depositor DCC
R_{free} test set	4579 reflections (7.61%)	DCC
Wilson B-factor (Å ²)	34.3	Xtriage
Anisotropy	0.619	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 40.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	5 of 64942 reflections (0.008%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10309	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 49.58 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 7.3114e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, ATP

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.88	0/5037	0.84	12/6818 (0.2%)
1	B	0.89	0/5021	0.83	9/6800 (0.1%)
All	All	0.89	0/10058	0.83	21/13618 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	6
1	B	0	5
All	All	0	11

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	502	ARG	NE-CZ-NH2	-10.43	115.08	120.30
1	B	502	ARG	NE-CZ-NH2	-9.35	115.63	120.30
1	A	502	ARG	NE-CZ-NH1	8.82	124.71	120.30
1	B	502	ARG	NE-CZ-NH1	7.53	124.06	120.30
1	A	500	LEU	CA-CB-CG	-7.08	99.01	115.30
1	A	861	ARG	NE-CZ-NH2	-7.04	116.78	120.30
1	A	297	ASP	N-CA-C	6.66	128.98	111.00
1	B	500	LEU	CA-CB-CG	-6.46	100.44	115.30
1	A	322	LEU	CA-CB-CG	6.23	129.63	115.30
1	A	411	LEU	CA-CB-CG	6.09	129.30	115.30
1	A	662	ARG	NE-CZ-NH2	-6.05	117.27	120.30
1	B	861	ARG	NE-CZ-NH1	6.05	123.33	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	707	ASP	CB-CG-OD1	5.97	123.68	118.30
1	B	778	ARG	NE-CZ-NH2	-5.91	117.35	120.30
1	A	660	VAL	CG1-CB-CG2	5.90	120.34	110.90
1	B	778	ARG	NE-CZ-NH1	5.87	123.23	120.30
1	B	744	ARG	NE-CZ-NH2	-5.61	117.49	120.30
1	B	662	ARG	NE-CZ-NH1	5.54	123.07	120.30
1	A	734	ARG	NE-CZ-NH1	5.52	123.06	120.30
1	B	662	ARG	NE-CZ-NH2	-5.28	117.66	120.30
1	A	861	ARG	NE-CZ-NH1	5.24	122.92	120.30

There are no chirality outliers.

All (11) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	408	ASN	Peptide
1	A	414	VAL	Peptide
1	A	415	PHE	Peptide
1	A	416	ALA	Peptide
1	A	417	GLU	Peptide
1	A	716	ASN	Peptide
1	B	408	ASN	Peptide
1	B	414	VAL	Peptide
1	B	415	PHE	Peptide
1	B	416	ALA	Peptide
1	B	417	GLU	Peptide

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4920	0	4819	83	0
1	B	4904	0	4786	74	0
2	A	31	0	12	4	0
2	B	31	0	12	5	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	200	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	219	0	0	3	0
All	All	10309	0	9629	160	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

All (160) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:1001:ATP:H4'	2:A:1001:ATP:O1B	1.52	1.10
1:B:531:LYS:HG2	1:B:706:HIS:HD2	1.22	0.99
1:A:602:ARG:HG3	1:A:604:SER:H	1.29	0.98
1:B:531:LYS:HG2	1:B:706:HIS:CD2	2.04	0.92
1:B:602:ARG:HG3	1:B:604:SER:H	1.33	0.90
1:B:782:ASN:HB3	1:B:888:MET:CE	2.06	0.86
1:A:782:ASN:HB3	1:A:888:MET:CE	2.08	0.82
1:A:469:LYS:HD2	1:A:712:PRO:CG	2.09	0.81
2:B:1001:ATP:H4'	2:B:1001:ATP:O1B	1.81	0.80
1:A:449:CYS:SG	1:A:452:CYS:HB2	2.21	0.80
1:B:752:ASN:HB2	1:B:755:ASP:H	1.48	0.79
1:A:653:PHE:C	1:A:653:PHE:CD2	2.58	0.77
1:B:449:CYS:SG	1:B:452:CYS:HB2	2.25	0.77
1:A:600:ASP:CG	1:A:601:GLN:H	1.90	0.76
1:B:322:LEU:HB2	1:B:746:SER:O	1.86	0.76
1:A:297:ASP:HB2	1:A:307:TYR:CE2	2.20	0.75
1:B:292:THR:HG22	1:B:293:THR:HG23	1.68	0.75
1:B:600:ASP:CG	1:B:601:GLN:H	1.90	0.74
2:A:1001:ATP:C4'	2:A:1001:ATP:O1B	2.33	0.74
1:B:723:MET:HG2	1:B:842:PRO:HG3	1.67	0.74
1:A:653:PHE:C	1:A:653:PHE:HD2	1.90	0.74
1:A:275:ASN:HD22	1:A:577:HIS:CD2	2.07	0.72
1:B:734:ARG:NH1	2:B:1001:ATP:O2G	2.21	0.72
1:B:278:LYS:HD2	1:B:577:HIS:ND1	2.04	0.71
1:A:469:LYS:HD2	1:A:712:PRO:HG2	1.71	0.71
1:A:666:SER:OG	1:A:803:HIS:HE1	1.74	0.70
1:A:506:GLY:HA3	1:A:660:VAL:HG13	1.74	0.69
1:B:469:LYS:HD2	1:B:712:PRO:CG	2.24	0.67
1:B:600:ASP:C	1:B:601:GLN:HG3	2.14	0.67
1:A:782:ASN:HB3	1:A:888:MET:HE1	1.75	0.66
1:A:522:ARG:HD2	4:A:1107:HOH:O	1.96	0.66
1:A:602:ARG:CG	1:A:603:GLY:N	2.60	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:602:ARG:HG3	1:B:603:GLY:N	2.12	0.64
1:B:600:ASP:O	1:B:601:GLN:HG3	1.97	0.64
1:B:782:ASN:CB	1:B:888:MET:HE1	2.28	0.64
1:A:531:LYS:HE2	1:A:706:HIS:CD2	2.33	0.64
1:B:653:PHE:C	1:B:653:PHE:CD2	2.70	0.64
1:A:292:THR:HG22	1:A:293:THR:HG23	1.81	0.63
1:A:297:ASP:OD2	1:A:300:HIS:N	2.32	0.62
1:B:604:SER:HA	1:B:609:THR:OG1	1.99	0.62
1:B:577:HIS:CD2	1:B:599:GLU:HG3	2.34	0.62
1:A:602:ARG:HG3	1:A:603:GLY:N	2.13	0.62
1:B:417:GLU:HG3	1:B:419:ASN:HB2	1.81	0.62
1:A:667:GLY:HA3	2:A:1001:ATP:HN61	1.65	0.62
1:A:322:LEU:HB2	1:A:746:SER:O	1.99	0.62
1:B:782:ASN:HB3	1:B:888:MET:HE1	1.81	0.61
1:A:782:ASN:CB	1:A:888:MET:HE1	2.31	0.61
1:A:278:LYS:HD2	1:A:577:HIS:ND1	2.16	0.60
1:A:653:PHE:HD2	1:A:654:GLU:N	1.99	0.60
1:B:602:ARG:CG	1:B:603:GLY:N	2.65	0.59
1:A:416:ALA:HB1	1:A:418:GLN:OE1	2.02	0.59
1:A:414:VAL:CG1	1:A:478:PHE:CE2	2.85	0.59
1:B:506:GLY:HA3	1:B:660:VAL:HG13	1.84	0.59
1:B:734:ARG:HH21	1:B:738:GLU:CD	2.06	0.58
1:B:416:ALA:H	1:B:478:PHE:HD1	1.49	0.58
1:B:653:PHE:O	1:B:653:PHE:CD2	2.56	0.58
1:A:404:LYS:HE3	4:A:1254:HOH:O	2.03	0.58
1:A:752:ASN:HB2	1:A:755:ASP:H	1.68	0.58
1:A:506:GLY:HA3	1:A:660:VAL:CG1	2.33	0.57
1:B:414:VAL:HG11	1:B:478:PHE:CE2	2.39	0.57
1:A:487:GLU:OE2	1:A:575:TYR:OH	2.23	0.57
1:A:602:ARG:CG	1:A:603:GLY:H	2.18	0.56
1:B:440:GLU:HA	1:B:443:ASN:HD22	1.69	0.56
1:A:742:ARG:HD2	4:A:1217:HOH:O	2.05	0.56
1:A:414:VAL:HG11	1:A:478:PHE:CE2	2.40	0.56
1:B:407:SER:O	1:B:409:ALA:N	2.39	0.56
1:B:322:LEU:N	1:B:322:LEU:HD13	2.20	0.55
1:A:734:ARG:HH21	1:A:738:GLU:CD	2.10	0.55
1:B:653:PHE:HD2	1:B:653:PHE:O	1.90	0.54
1:A:604:SER:HA	1:A:609:THR:OG1	2.08	0.54
1:A:772:PHE:O	1:A:778:ARG:HD3	2.08	0.53
1:A:666:SER:OG	1:A:803:HIS:CE1	2.61	0.53
1:B:506:GLY:HA3	1:B:660:VAL:CG1	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:761:LYS:HG2	1:A:794:PRO:HG3	1.91	0.53
1:B:782:ASN:CB	1:B:888:MET:CE	2.82	0.53
1:B:602:ARG:CG	1:B:603:GLY:H	2.21	0.53
1:A:298:PRO:HD2	1:A:584:ARG:HH22	1.75	0.52
1:A:600:ASP:CG	1:A:601:GLN:N	2.61	0.52
1:A:723:MET:HG2	1:A:842:PRO:HG3	1.92	0.51
1:A:487:GLU:OE2	1:A:607:VAL:N	2.43	0.51
1:A:600:ASP:O	1:A:601:GLN:HG3	2.10	0.51
1:B:414:VAL:CG1	1:B:478:PHE:CE2	2.93	0.51
1:A:477:TRP:HB2	1:A:605:GLY:HA3	1.93	0.51
1:B:531:LYS:CG	1:B:706:HIS:CD2	2.88	0.51
1:A:544:ILE:HD11	1:A:602:ARG:HB3	1.94	0.50
1:A:349:THR:O	1:A:353:GLN:HG3	2.11	0.50
1:B:487:GLU:OE2	1:B:575:TYR:OH	2.28	0.50
1:B:600:ASP:CG	1:B:601:GLN:N	2.62	0.50
1:B:469:LYS:HD2	1:B:712:PRO:HG2	1.93	0.50
1:B:522:ARG:HD2	4:B:1115:HOH:O	2.11	0.50
1:A:297:ASP:HB2	1:A:307:TYR:CZ	2.47	0.50
1:B:554:LYS:O	1:B:557:GLU:HG2	2.11	0.50
1:B:734:ARG:NH2	2:B:1001:ATP:O2G	2.44	0.49
1:A:315:ALA:HB2	1:A:591:THR:OG1	2.12	0.49
1:B:577:HIS:HD2	1:B:599:GLU:HG3	1.74	0.49
1:A:782:ASN:HB3	1:A:888:MET:HE3	1.92	0.48
1:A:415:PHE:HB2	1:A:416:ALA:HB2	1.95	0.48
1:B:462:LYS:HG3	4:B:1171:HOH:O	2.14	0.48
1:A:544:ILE:HD11	1:A:602:ARG:CB	2.44	0.48
1:A:443:ASN:ND2	1:A:446:ARG:HH12	2.12	0.48
1:B:602:ARG:HD2	1:B:609:THR:HG23	1.95	0.48
1:B:668:ASP:H	2:B:1001:ATP:HN61	1.61	0.47
1:B:460:ARG:NH2	1:B:463:LYS:HD2	2.30	0.47
1:B:407:SER:C	1:B:409:ALA:H	2.18	0.47
1:A:393:PRO:HG2	1:A:558:LEU:HB3	1.97	0.47
1:B:468:GLY:O	1:B:536:ASP:HB2	2.15	0.47
1:B:463:LYS:HE3	2:B:1001:ATP:O3A	2.15	0.47
1:B:396:CYS:HB3	1:B:488:PHE:CE2	2.51	0.46
1:B:477:TRP:HB2	1:B:605:GLY:HA3	1.97	0.46
1:B:416:ALA:HB1	1:B:418:GLN:OE1	2.16	0.45
1:A:298:PRO:HB2	1:A:584:ARG:HH12	1.81	0.45
1:B:407:SER:C	1:B:409:ALA:N	2.70	0.45
1:A:416:ALA:CB	1:A:418:GLN:OE1	2.64	0.45
1:B:546:ARG:O	1:B:550:GLU:HG3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:414:VAL:CG1	1:B:478:PHE:CZ	3.00	0.45
1:B:782:ASN:HB3	1:B:888:MET:HE3	1.94	0.44
1:A:443:ASN:HD22	1:A:446:ARG:NH1	2.15	0.44
1:B:487:GLU:OE2	1:B:607:VAL:HB	2.17	0.44
1:A:443:ASN:HD22	1:A:446:ARG:HH12	1.64	0.44
1:A:287:LYS:HG3	1:A:294:TRP:CE2	2.53	0.44
1:A:557:GLU:HG2	1:A:557:GLU:H	1.52	0.44
1:A:600:ASP:C	1:A:601:GLN:HG3	2.38	0.44
1:A:820:ARG:HA	1:A:824:GLU:HB3	2.00	0.43
1:B:604:SER:CB	1:B:609:THR:HG21	2.48	0.43
1:A:416:ALA:H	1:A:478:PHE:HD1	1.64	0.43
1:A:603:GLY:O	1:A:604:SER:HB3	2.18	0.43
1:A:463:LYS:HE3	2:A:1001:ATP:O3A	2.18	0.43
1:A:468:GLY:O	1:A:536:ASP:HB2	2.19	0.42
1:B:487:GLU:OE2	1:B:607:VAL:N	2.52	0.42
1:B:457:MET:HB2	1:B:478:PHE:CD2	2.54	0.42
1:A:598:ARG:HG2	1:A:600:ASP:HB2	2.00	0.42
1:B:617:ASN:O	1:B:621:GLN:HG2	2.18	0.42
1:A:519:TYR:HB3	1:A:828:TRP:CE2	2.54	0.42
1:A:425:ARG:HB2	1:A:425:ARG:HE	1.68	0.42
1:A:294:TRP:HA	1:A:310:SER:O	2.20	0.42
1:B:742:ARG:HD2	4:B:1142:HOH:O	2.20	0.42
1:B:773:HIS:CD2	1:B:773:HIS:H	2.38	0.42
1:B:604:SER:HA	1:B:609:THR:HG21	2.01	0.42
1:B:519:TYR:HB3	1:B:828:TRP:CE2	2.54	0.42
1:A:773:HIS:CD2	1:A:773:HIS:H	2.37	0.42
1:A:353:GLN:HA	1:A:456:MET:HE1	2.01	0.42
1:B:557:GLU:H	1:B:557:GLU:HG2	1.60	0.41
1:B:332:LYS:HB3	1:B:333:PRO:HD3	2.02	0.41
1:A:787:ALA:HB1	1:A:875:VAL:HB	2.02	0.41
1:A:531:LYS:HE2	1:A:706:HIS:HD2	1.79	0.41
1:A:418:GLN:NE2	4:A:1296:HOH:O	2.52	0.41
1:B:558:LEU:N	1:B:558:LEU:HD22	2.36	0.41
1:A:457:MET:HB2	1:A:478:PHE:CD2	2.56	0.41
1:A:409:ALA:C	1:A:411:LEU:N	2.75	0.41
1:B:500:LEU:HD22	1:B:614:THR:CG2	2.51	0.41
1:A:469:LYS:HD2	1:A:712:PRO:HG3	1.96	0.41
1:B:604:SER:HB2	1:B:609:THR:HG21	2.02	0.40
1:A:869:TYR:OH	1:A:888:MET:HB2	2.22	0.40
1:A:298:PRO:O	1:A:299:GLU:C	2.59	0.40
1:A:454:TYR:CE2	1:A:578:LYS:HG2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:402:ILE:HG12	1:A:428:VAL:HB	2.03	0.40
1:B:468:GLY:O	1:B:469:LYS:HE2	2.21	0.40
1:A:454:TYR:HB2	1:A:580:VAL:HG22	2.04	0.40
1:A:388:SER:CB	1:A:558:LEU:HD11	2.52	0.40
1:B:278:LYS:HD2	1:B:577:HIS:CE1	2.55	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	609/639 (95%)	572 (94%)	30 (5%)	7 (1%)	17	17
1	B	609/639 (95%)	575 (94%)	29 (5%)	5 (1%)	24	26
All	All	1218/1278 (95%)	1147 (94%)	59 (5%)	12 (1%)	19	20

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	420	GLN
1	B	408	ASN
1	B	411	LEU
1	A	297	ASP
1	A	408	ASN
1	A	411	LEU
1	B	409	ALA
1	B	420	GLN
1	B	422	SER
1	A	409	ALA
1	A	422	SER
1	A	587	ALA

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	515/544 (95%)	483 (94%)	32 (6%)	23	28
1	B	512/544 (94%)	478 (93%)	34 (7%)	21	25
All	All	1027/1088 (94%)	961 (94%)	66 (6%)	22	26

All (66) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	299	GLU
1	A	313	VAL
1	A	322	LEU
1	A	355	ARG
1	A	394	ARG
1	A	407	SER
1	A	414	VAL
1	A	415	PHE
1	A	432	ARG
1	A	481	LEU
1	A	521	LEU
1	A	528	GLN
1	A	546	ARG
1	A	557	GLU
1	A	558	LEU
1	A	588	GLU
1	A	590	LYS
1	A	601	GLN
1	A	602	ARG
1	A	635	GLN
1	A	653	PHE
1	A	660	VAL
1	A	753	VAL
1	A	778	ARG
1	A	804	SER
1	A	835	ILE
1	A	839	THR

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Mol	Chain	Res	Type
1	A	847	ARG
1	A	860	SER
1	A	861	ARG
1	A	874	GLN
1	A	888	MET
1	B	285	LYS
1	B	299	GLU
1	B	313	VAL
1	B	322	LEU
1	B	332	LYS
1	B	394	ARG
1	B	407	SER
1	B	414	VAL
1	B	415	PHE
1	B	426	GLU
1	B	432	ARG
1	B	457	MET
1	B	481	LEU
1	B	521	LEU
1	B	544	ILE
1	B	557	GLU
1	B	588	GLU
1	B	590	LYS
1	B	601	GLN
1	B	602	ARG
1	B	635	GLN
1	B	653	PHE
1	B	660	VAL
1	B	778	ARG
1	B	793	VAL
1	B	804	SER
1	B	835	ILE
1	B	839	THR
1	B	847	ARG
1	B	848	GLU
1	B	860	SER
1	B	861	ARG
1	B	874	GLN
1	B	888	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	443	ASN
1	A	577	HIS
1	A	706	HIS
1	A	803	HIS
1	B	706	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	ATP	A	1001	-	24,33,33	1.80	2 (8%)	31,52,52	3.26	9 (29%)
2	ATP	B	1001	-	24,33,33	1.78	4 (16%)	31,52,52	3.05	8 (25%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ATP	A	1001	-	-	0/18/38/38	0/3/3/3
2	ATP	B	1001	-	-	0/18/38/38	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1001	ATP	C5'-C4'	2.32	1.59	1.51
2	B	1001	ATP	PG-O3G	2.44	1.63	1.54
2	B	1001	ATP	PG-O2G	3.09	1.65	1.54
2	A	1001	ATP	PG-O3G	3.27	1.66	1.54
2	B	1001	ATP	O4'-C1'	6.48	1.49	1.41
2	A	1001	ATP	O4'-C1'	7.15	1.50	1.41

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1001	ATP	N3-C2-N1	-12.08	119.64	128.89
2	A	1001	ATP	N3-C2-N1	-11.55	120.05	128.89
2	A	1001	ATP	PB-O3B-PG	-8.60	103.84	132.67
2	B	1001	ATP	PB-O3B-PG	-6.13	112.13	132.67
2	B	1001	ATP	PA-O3A-PB	-5.44	117.45	132.73
2	A	1001	ATP	PA-O3A-PB	-5.42	117.50	132.73
2	B	1001	ATP	C4-C5-N7	-4.49	105.35	109.48
2	A	1001	ATP	C4-C5-N7	-4.17	105.64	109.48
2	A	1001	ATP	C5'-C4'-C3'	-3.33	102.00	115.21
2	A	1001	ATP	C2'-C1'-N9	-3.25	109.32	114.29
2	B	1001	ATP	C1'-N9-C4	-2.55	123.09	126.94
2	A	1001	ATP	C1'-N9-C4	-2.08	123.80	126.94
2	A	1001	ATP	C2-N1-C6	2.13	122.58	118.77
2	B	1001	ATP	O2B-PB-O1B	2.14	124.14	112.53
2	B	1001	ATP	C2-N1-C6	2.20	122.70	118.77
2	B	1001	ATP	O2B-PB-O3A	2.38	115.89	105.09
2	A	1001	ATP	O4'-C1'-N9	3.99	116.45	108.10

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1001	ATP	4	0
2	B	1001	ATP	5	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	613/639 (95%)	0.09	31 (5%) 32 39	25, 45, 74, 94	0
1	B	613/639 (95%)	0.04	24 (3%) 43 51	26, 46, 73, 92	0
All	All	1226/1278 (95%)	0.07	55 (4%) 37 45	25, 45, 73, 94	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	419	ASN	6.0
1	A	748	GLY	5.9
1	A	416	ALA	5.3
1	B	416	ALA	5.1
1	B	890	SER	5.1
1	B	751	TRP	4.8
1	A	751	TRP	4.7
1	A	413	ALA	4.5
1	B	419	ASN	4.3
1	A	750	GLY	4.2
1	B	415	PHE	3.9
1	A	587	ALA	3.9
1	A	415	PHE	3.8
1	B	414	VAL	3.6
1	B	830	MET	3.6
1	A	299	GLU	3.6
1	A	315	ALA	3.4
1	B	411	LEU	3.4
1	A	412	GLY	3.3
1	A	414	VAL	3.2
1	A	603	GLY	3.1
1	A	749	ALA	3.1
1	B	421	TRP	3.1
1	B	889	THR	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	298	PRO	3.0
1	A	420	GLN	3.0
1	A	888	MET	2.9
1	A	286	LEU	2.9
1	A	422	SER	2.8
1	B	315	ALA	2.8
1	A	466	GLU	2.8
1	B	750	GLY	2.7
1	B	418	GLN	2.6
1	B	410	ALA	2.6
1	B	413	ALA	2.6
1	B	420	GLN	2.6
1	A	418	GLN	2.5
1	A	869	TYR	2.5
1	B	466	GLU	2.4
1	A	421	TRP	2.4
1	B	706	HIS	2.3
1	B	275	ASN	2.3
1	B	749	ALA	2.2
1	B	603	GLY	2.1
1	B	851	TRP	2.1
1	A	830	MET	2.1
1	B	590	LYS	2.1
1	A	851	TRP	2.1
1	A	411	LEU	2.1
1	A	588	GLU	2.1
1	A	889	THR	2.1
1	A	890	SER	2.1
1	A	590	LYS	2.1
1	A	478	PHE	2.0
1	B	286	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	ATP	B	1001	31/31	0.83	0.23	4.84	48,63,68,71	0
2	ATP	A	1001	31/31	0.84	0.21	3.11	58,68,73,74	0
3	ZN	B	1003	1/1	0.99	0.02	-1.78	39,39,39,39	0
3	ZN	A	1003	1/1	0.99	0.01	-3.11	39,39,39,39	0
3	ZN	B	1002	1/1	0.99	0.02	-5.13	33,33,33,33	0
3	ZN	A	1002	1/1	0.99	0.02	-7.19	28,28,28,28	0

6.5 Other polymers [i](#)

There are no such residues in this entry.